

<p><b>96-287086/29</b> C03 (C02) <b>FARB 94.12.08</b>  <b>BAYER AG</b> *WO 9617825-A1  94.12.08 94DE-4443641 (96.06.13) C07C 251/48, A01N 37/50, 43/40, 43/54, C07C 255/44, C07D 285/08, 417/04, 239/34, C07C 251/60, A01N 43/82, C07C 69/734  <b>New aromatic and araliphatic carboxamide cpds. - useful as fungicides in plant protection, esp. for control of cereal, fruit and vegetable diseases (Ger)</b>  <b>C96-091791</b> N(AU BB BG BR BY CA CN CZ FI HU JP KR KZ LK MX NO NZ PL RO RU SK UA US) R(AT BE CH DE DK ES FR GB GR IE IT LU MC NL OA PT SE)  Addnl. Data: SEITZ T, HEINEMANN U, STENZEL K, DUTZMANN S  95.11.27 95WO-EP04668</p>	<p>C(7-H, 10-D3, 14-A6) .3</p> <p>E = 2-R<sup>1</sup>-1-alkene-1,1-diyl, 2-aza-2-R<sup>2</sup>-1-alkene-1,1-diyl, NR<sup>3</sup>, 3-aza-1-R<sup>4</sup>-3-R<sup>5</sup>-1-propene-2,3-diyl, 3-(aza or thia)-1-R<sup>4</sup>-1-propenediyl, 1-aza-1-(R<sup>4</sup> or R<sup>6</sup>)-3-R<sup>5</sup>-1-propene-2,3-diyl, 1,3-diaza-1,3-R<sup>5</sup>-1-propene, 2,3-diyl, or 1-aza-3-(oxa or thia)-1-R<sup>6</sup>-1-propene-2,3-diyl;  R<sup>1</sup>, R<sup>4</sup> = alkyl, alkoxy, alkylthio, alkylamino or dialkylamino (all opt. substd.), H, halo or CN;  R<sup>2</sup>, R<sup>6</sup> = alkyl, alkoxy, alkylamino or dialkylamino (all opt. substd.), H, NH<sub>2</sub> or CN;  R<sup>3</sup> = alkyl, alkenyl, alkynyl, cycloalkyl or cycloalkylalkyl (all opt. substd.), H or CN;  R<sup>5</sup> = alkyl;  G = alkanediyl, alkenediyl or alkynediyl (all opt. substd. by halo, OH, haloalkyl or cycloalkyl), bond, O, S, -Q-CQ-, -CQ-Q-, -CH<sub>2</sub>-Q-, -Q-CH<sub>2</sub>-, -CQ-Q-CH<sub>2</sub>-, -CH<sub>2</sub>-Q-CQ-; -Q-CQ-CH<sub>2</sub>-, -Q-CQ-Q-CH<sub>2</sub>-, -N=N-, -S(O)<sub>n</sub>-, -CH<sub>2</sub>-S(O)<sub>n</sub>-, -CQ-, -S(O)<sub>n</sub>-CH<sub>2</sub>-, -C(R<sup>7</sup>)=N-O-, -C(R<sup>7</sup>)=N-O-CH<sub>2</sub>-, -N(R<sup>8</sup>)-, -CQ-N(R<sup>8</sup>)-, -N(R<sup>8</sup>)-CQ-, -Q-CQ-, -N(R<sup>8</sup>)-CQ-Q-, -N(R<sup>8</sup>)-CQ-Q-, -CH<sub>2</sub>-O-N=C(R<sup>7</sup>)-, -N(R<sup>8</sup>)-CQ-Q-, -CQ-N(R<sup>8</sup>)-CQ-Q-, -N(R<sup>8</sup>)-CQ-Q-CH<sub>2</sub>-, -Q-C(R<sup>7</sup>)=N-O-CH<sub>2</sub>- or -</p> <p>WO 9617825-A+</p>
<p>Substd (hetero)aryl and (hetero)aralkyl carboxamides of formula (I) are new.</p> <p style="text-align: center;">Z-G-Ar<sup>1</sup>-E-CO-N(A<sup>1</sup>)-[C(A<sup>2</sup>)(A<sup>3</sup>)]<sub>m</sub>-Ar<sup>2</sup> (I)</p> <p>A<sup>1</sup>, A<sup>2</sup> = H or alkyl;  A<sup>3</sup> = H, alkyl or CN;  Ar<sup>1</sup> = opt. substd. arylene or heteroarylene;  Ar<sup>2</sup> = opt. substd. aryl or heteroaryl;</p>	

<p> <math>N(R^8)-C(R^7)=N-O-CH_2-</math>;  <math>m, n = 0-2</math>;  <math>Q = O</math> or <math>S</math>;  <math>R^7 =</math> alkyl, alkoxy, alkylthio, alkylamino, dialkylamino or cycloalkyl  (all opt. substd.), <math>H</math> or <math>CN</math>;  <math>R^8 =</math> alkyl, alkoxy or cycloalkyl (all opt., substd.), <math>H</math>, <math>OH</math> or <math>CN</math>;  <math>Z =</math> alkyl, alkanyl, alkynyl, cycloalkyl, aryl or heterocyclyl (all opt.  substd.). </p>	<p> <math>S(O)_{n-}, -CQ-, -S(O)_n-CH_2-, -C(R^7)=N-O-, -C(R^7)=N-O-CH_2-, -</math>  <math>N(R^8)-, -CQ-N(R^8)-, -N(R^8)-CQ-, -Q-CQ-N(R^8)-, -N=C(R^7)-Q-</math>  <math>CH_2-, -CH_2-O-N=C(R^7)-, -N(R^8)-CQ-Q-, -CQ-N(R^8)-CQ-Q-, -</math>  <math>N(R^8)-CQ-Q-CH_2-, -Q-C(R^7)=N-O-CH_2-</math> or <math>-N(R^8)-C(R^7)=N-O-</math>  <math>CH_2-</math>;  <math>R^7 =</math> <math>Cy, T'', OT'', ST'', NHT'', NT''_2, H</math>, halo or <math>CN</math>;  <math>R^8 =</math> <math>Cy, T'', H, OH</math> or <math>CN</math>;  <math>Z =</math> 1-8C alkyl (opt. substd. by 1 or more halo, <math>CN, OH, NH_2, OT''</math>,  <math>ST'</math>, <math>SOT'</math> or <math>SO_2T'</math> (opt. substd. by halo)), 2-8C alkenyl (opt.  substd. by halo), 2-8C alkynyl (opt. substd. by halo), 3-6C  cycloalkyl (opt. substd. by one or more halo, <math>CN, COOH, Ph, T'</math>,  <math>COOT''</math>), or <math>Ar</math>;  <math>Ph =</math> phenyl (opt. substd. by halo, <math>CN, T'</math>, 1-4C haloalkyl, 1-4C  haloalkoxy or <math>OT''</math>);  <math>Ar =</math> phenyl, naphthyl or 3-7-membered heterocycl contg. <math>O, S</math> or <math>N</math>  and opt. 1 or 2 additional <math>N</math> (all opt. substd. by one or more <math>R^9</math>);  <math>R^0 =</math> halo, <math>CN, NO_2, NH_2, OH, CHO, COOH, CONH_2, CSNH_2, XR_n, 2-</math>  6C alkenyl, 2-6C alkenyloxy (both opt. substd. by 1-13 halo),  <math>NHT, NT''_2, COT, OCOT, COOT, OSO_2T, T</math> (substd. by <math>NH_2OH</math>  or <math>NHOT</math>), <math>Cy, Alk, OAlkO, Het, CH_2Het</math> or <math>Ar^3</math>;  <math>Alk =</math> 1-6C alkylene (opt. substd. by one or more halo, or <math>T'</math> (opt. </p>
<p> <b>MORE SPECIFICALLY</b>  <math>R^1, R^4 = T'', OT'', ST'', NHT'', NT''_2, H</math>, halo or <math>CN</math>;  <math>R^2, R^6 = T'', OT'', ST'', NHT'', NT''_2, H</math>, halo or <math>CN</math>;  <math>R^3 =</math> 2-6C alkenyl, 2-6C alkynyl (both opt. substd. as for <math>T''</math>), <math>T'', Cy-</math>  substd. 1-4C alkyl, <math>H</math> or <math>CN</math>;  <math>T'' = T</math> (opt. substd. by halo, <math>CN</math> or <math>OT''</math>)  <math>T, R^5 =</math> 1-6C alkyl;  <math>T' =</math> 1-4C alkyl;  <math>Cy =</math> 3-6C cycloalkyl (opt. substd. by halo, <math>CN, COOH, T'</math> or <math>COOT'</math>);  <math>G =</math> 1-4C alkanediyl, 2-4C alkenediyl or 2-4C alkynediyl (all opt.  substd. by halo, <math>OH, T', 1-4C</math> haloalkyl or 3-6C cycloalkyl),  bond, <math>O, S, -Q-CQ-, -CQ-Q-, -CH_2-Q-, -Q-CH_2-, -CQ-Q-CH_2-, -</math>  <math>CH_2-Q-CQ-, -Q-CQ-CH_2-, -Q-CQ-Q-CH_2-, -N=N-, -S(O)_n-, -CH_2-</math> </p>	

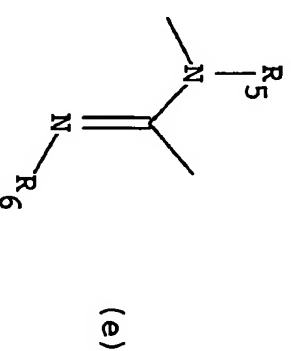
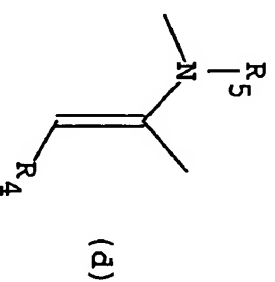
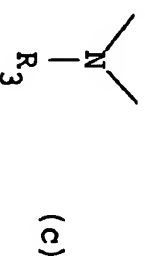
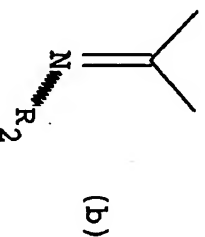
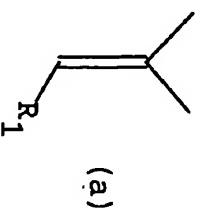
subst. by 1-9 halo));

X = bond, O, S, SO, SO<sub>2</sub>;

Het = 3-7 membered heterocycle with 1-3 heteroatoms (esp. N, O S);  
Ar<sup>3</sup> = pyridyl, thienyl, phenyl, phenoxy, phenylthio, benzyl,  
benzyloxy, benzylthio, phenylethyl or phenylethyloxy (all opt.  
ring-subst. by R");

R" = T', ST', OT' (all opt. subst. with 1-9 halo), halo, CN, Alk or  
OAlkO;

E = a gp. of formula (a)-(e).



### USE

(1) are plant fungicides which are tolerated well by plants and which are esp. useful for the control of cereal diseases caused by various spp. including *Erysiphe*, *Leptosphaeria*, *Pyrenophora* and *Cochliobolus* spp. as well as diseases in fruit and vegetable crops caused by various spp. including *Podosphaera*.

Application rate is 0.001-50g/kg when used as a seed dressing.

### ADVANTAGE

(I) are more effective than known substd. carboxamides (cf. e.g. EP 398692) and have esp. good activity in vitro.

### PREPARATION

Claimed prepn. of (I) is as follows.



(II)

(III)

R = OH, halo or alkoxy.

(III) can be used as a hydrohalide and the reaction may be performed in the presence of an acid acceptor, condensation agent and/or diluent.

### EXAMPLE

A mixt. of methyl 2-methoxyimino-2-[2-(2-methylphenoxy-methyl)-phenyl]-acetate (2.5 g) and 4-chlorobenzylamine (1.14 g) was stirred at 120 °C for 12 hrs., cooled, taken up in CH<sub>2</sub>Cl<sub>2</sub>, washed (H<sub>2</sub>O, 1N HCl, then H<sub>2</sub>O), dried (Na<sub>2</sub>SO<sub>4</sub>) and filtered. Concentration and chromatography (SiO<sub>2</sub>; petroleum ether:EtOAc, 5:1) gave N-(4-chlorobenzyl)-2-methoxyimino-2-[2-(2-methylphenoxy)methyl)-

acetamide (1.4 g; 42% yield); oil.

Typical cpds. (I) applied at 250 g/ha gave 100% protection to wheat and barley from attack by Erysiphe graminis and also gave 100% kill of this fungus on the same cereals. When used in concns. of 20 ppm, cpds. also gave 89-97% protection to apples from attack by Podosphaera leucotricha. (LJ)  
(83pp2101DwgNo. 0/0)  
SR:6.Jnl.Ref EP431328 JP01031753 JP02142761 WO9501328